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Sulphur, molecular complexity of (KELLAS), 903.

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Tetrahydronaphthalene series (GREEN and ROWE), 955.
Thermotropy and phototropy (SENIER and GALLAGHER), 28.
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FORMULA INDEX.

THE following index of organic compounds of known empirical formula is arranged according to Richter's system (see *Lexikon der Kohlenstoff-Verbindungen*).

The elements are given in the order C, H, O, N, Cl, Br, I, F, S, P, and the remainder alphabetically.

The compounds are arranged—

Firstly, in groups according to the number of carbon atoms (thus C_1 group, C_2 group, etc.).

Secondly, according to the number of other elements besides carbon contained in the molecule (thus 5 IV indicates that the molecule contains five carbon atoms and four other elements).

Thirdly, according to the nature of the elements present in the molecule (given in the above order).

Fourthly, according to the number of atoms of each single element (except carbon) present in the molecule.

Salts are placed with the compounds from which they are derived. The chlorides, bromides, iodides, and cyanides of quaternary ammonium bases, however, are registered as group-substances.

C_1 Group.

CH_4 Methane, ignition of mixtures of air and (MASON and WHEELER), 45; (WHEELER), 840.

CCl_4 Carbon tetrachloride, action of aniline with (HARTUNG), 163.

1 II

CH_2O Formaldehyde, action of carbamide and (DIXON), 238.

CH_3I Methyl iodide, relative activities of ethyl iodide, propyl iodide and, with sodium α- and β-naphthoxides (COX), 666.

$COCl_2$ Carbonyl chloride, action of ammonia with (WERNER), 694.

1 III

CHON Cyanic acid, metallic salts, hydrolysis of, 84.

CH_4ON_2 Carbamide, synthesis of (WERNER and CARPENTER), 694; mechanism of the synthesis of, from urethane (WERNER), 622; decomposition of, on heating with acids and with alkalis (WERNER), 84; interaction of formaldehyde and (DIXON), 238.

C_2 Group.

C_2H_5I Ethyl iodide, relative activities of methyl iodide, propyl iodide and, with sodium α- and β-naphthoxides (COX), 666.

C_2H_6O Ethyl alcohol, association of organic compounds in solution in (INNES), 410.

C_2H_7N Ethylamine, preparation of (WERNER), 899.

2 III

$C_2H_4O_2N_2$ Oxamide, formation of urea and of biuret from (WERNER and CARPENTER), 693.

$C_2H_6ON_2$ Glycollamidine, salts of (RULE), 17.

$C_2H_6O_2N_2$ Methylolcarbamide, preparation of (DIXON), 246.

C_3 Group.

C_3H_7I *n*-Propyl iodide, relative activities of methyl iodide, ethyl iodide and with sodium α- and β-naphthoxides (COX), 666.

3 III

$C_3H_7O_2N$ Urethane, mechanism of the synthesis of urea from (WERNER), 622.

C_4 Group.

C_4H_9Cl *n*-Butyl chloride, action of, with *o*- and *p*-toluidine (REILLY and HICKINBOTTOM), 974.

$C_4H_{11}N$ Diethylamine, preparation of (WERNER), 899; stannochloride of (DRUCE), 715.

4 III

$C_4H_{10}O_2N_2$ Acetiminohydrin, preparation of (RULE), 11.

C_5 Group.

$C_5H_{10}O_3N_4$ Dimethylolcarbamide, preparation of (DIXON), 247.

$C_5H_{11}O_2N$ Methoxyacetiminoethyl ether (RULE), 9.

C_6 Group.

C_6H_6 Benzene, association of organic compounds in solution in (INNES), 410.

6 II

C_6H_6O Phenol, freezing points of mixtures of cresols and (DAWSON and MOUNTFORD), 923; estimation of, in mixtures with cresols (DAWSON and MOUNTFORD), 935.

C_6H_7N Aniline, action of carbon tetrachloride with (HARTUNG), 163.

$C_6H_8N_2$ *m*- and *p*-Phenylenediamine, stanni- and stanno-chlorides of (DRUCE), 716.

$C_6H_{10}O_4$ Dimethyl tartrate, association of, in benzene (INNES), 433.

$C_6H_{12}N_2$ Formaldehydepiperylhydrazone (WEINHAGEN), 586.

6 III

$C_6H_3O_6N_4$ 3:5-Dinitrobenz*iso*oxadiazole oxide (GREEN and ROWE), 70.

$C_6H_{14}O_4N_2$ Methoxyacetiminohydrin (RULE), 9.

C_7 Group.

$C_7H_6O_2$ Benzoic acid, calcium salt, hydrates and alcoholates of (STANBRIDGE), 805.

C_7H_8O *o*-, *m*-, and *p*-Cresol, freezing points of mixtures of phenol and (DAWSON and MOUNTFORD), 923; estimation of, in mixtures with phenol (DAWSON and MOUNTFORD), 935.

C_7H_9N Benzylamine, stanni- and stanno-chlorides of (DRUCE), 717.

Methylaniline, stanni- and stanno-chlorides of (DRUCE), 716.

o-Toluidine, stanni- and stanno-chlorides of (DRUCE), 716.

o- and *p*-Toluidine, action of *n*-butyl chloride on (REILLY and HICKINBOTTOM), 974.

$C_7H_{14}O_6$ β-Methylfructoside, preparation and structure of (STEELE), 257.

γ-Methylgalactoside (CUNNINGHAM), 598.

7 III

$C_7H_5O_8N_3$ Trinitroguaiacol, and its pyridine salt (POLLECOFF and ROBINSON), 653.

$C_7H_6O_6N_2$ Dinitroguaiacols (POLLECOFF and ROBINSON), 649.

C_7H_5ON Formanilide, association of, in benzene (INNES), 432.

$C_7H_8O_3S$ Toluene-*o*- and -*p*-sulphonic acids, analysis of mixtures of (MCKIE), 799.

$C_7H_9O_5N_4$ 3:5-Dinitro-2:6-diaminoanisole (POLLECOFF and ROBINSON), 655.

$C_7H_7O_2N$ Phenylaminoacetic acid, sign of the product in conversion of phenylbromoacetic acid into (SENTER and TUCKER), 140.

7 IV

$C_7H_9O_2NS$ Toluene-*o*- and -*p*-sulphonamides, freezing point curves of mixtures of (McKIE), 799.

C_8 Group.

$C_8H_6O_4$ Terephthalic acid, menthyl alkyl esters of (COHEN and DE PENNINGTON), 57.

$C_8H_8O_2$ Phenylacetic acid, association of, in benzene (INNES), 432.

$C_8H_8O_3$ Methyl salicylate, association of, in benzene (INNES), 431.

$C_8H_{11}N$ *p*-Methylbenzylamine, stannochloride of (DRUCE), 718.

8 III

$C_8H_5O_6N$ Nitroterephthalic acid, menthyl alkyl esters of (COHEN and DE PENNINGTON), 57.

$C_8H_6O_2N_4$ 4-Acetyl-2- and -3-*p*-phenylene-1-diazo-4-imides (MORGAN and CLEAGE), 591.

$C_8H_7O_2Br$ Phenylbromoacetic acid, sign of the product in conversion of, into phenylaminoacetic acid (SENTER and TUCKER), 140.

$C_8H_7O_5N$ 6-Nitro-2-methoxybenzoic acid (SIMONSEN), 782.

$C_8H_7O_8N_3$ 3:5:6-Trinitroveratrole (POLLECOFF and ROBINSON), 654.

$C_8H_9O_3N_3$ *p*-Nitromethylformanilide (MORGAN and GRIST), 690.

$C_8H_8O_6N_2$ 5:6-Dinitroguaiacol methyl ether (POLLECOFF and ROBINSON), 651.

C_8H_9ON Acetanilide, association of, in benzene (INNES), 433.

Acetophenoneoxime, association of, in benzene (INNES), 432.

$C_8H_9O_2N$ Mandeliminohydrin (MACKENZIE), 1; (RULE), 12.

$C_8H_9O_2N_3$ Formylmethylaminobenzene-4-diazohydroxide (MORGAN and GRIST), 692.

$C_8H_9O_2Br$ 3-Bromoveratrole (SIMONSEN and RAU), 785.

$C_8H_9O_3N$ 6-Nitro-*o*-tolyl methyl ether (SIMONSEN), 781.

$C_8H_{10}ON_2$ *p*-Aminomethylformanilide (MORGAN and GRIST), 691.

$C_8H_{10}O_4N_2$ 5-Nitro-4-aminoveratrole (SIMONSEN and RAU), 27.

8 IV

$C_8H_5ON_3Cl_2$ 4-Acetyldichloro-*p*-phenylene-1-diazo-4-imides (MORGAN and CLEAGE), 594.

$C_8H_6O_2NCl$ *p*-Chloromandelamide, and its salts (RULE), 17.

$C_8H_{10}O_2NBr$ Bromoaminoveratroles (SIMONSEN and RAU), 786.

C_9 Group.

$C_9H_8O_2$ Cinnamic acid, association of, in benzene (INNES), 431.

$C_9H_{10}O_3$ Ethyl *p*-hydroxybenzoate, association of, in benzene (INNES), 434.

2-Methoxy-*m*-toluic acid, and its silver salt (SIMONSEN), 779.

9 III

$C_9H_9O_2Br$ α-Bromo-β-phenylpropionic acid, sign of the product in conversion of, into α-amino-β-phenylpropionic acid (SENTER, DREW, and MARTIN), 151.

$C_9H_9O_4N$ 5-Nitro-2- and -6-methoxy-*m*-tolualdehydes (SIMONSEN), 778.

$C_9H_9O_5N$ 5-Nitro-2- and -6-methoxy-*m*-toluic acids, and their salts (SIMONSEN), 779.

$C_9H_{10}O_3N_2$ 4-Nitroaceto-*o*-toluidide, trimorphism of (CHATTAWAY), 897.

$C_9H_{10}O_6N_2$ 6-Nitro-5-amino-3:4-dimethoxybenzoic acid (SIMONSEN and RAU), 26.

$C_9H_{10}O_7N_2$ 4:6-Dinitro-1:2:3-trimethoxybenzene (POLLECOFF and ROBINSON), 656.

$C_{11}H_{11}O_2N$ α-Amino-β-phenylpropionic acid, sign of the product in conversion of α-bromo-β-phenylpropionic acid into (SENTER, DREW, and MARTIN), 151.
2-Methoxy-*m*-tolualdoxime (SIMONSEN), 777.
$C_{11}H_{11}O_2N_3$ Acetylmethylaminobenzene-4-diazohydroxide (MORGAN and GRIST), 692.
$C_{11}H_{15}O_4N$ 5-Amino-3:4-dimethoxybenzoic acid, and its salts (SIMONSEN and RAU), 24.
$C_{11}H_{17}ON_3$ *p*-Aminomethylacetanilide (MORGAN and GRIST), 691.
$C_{11}H_{12}O_7N_4$ Nitrosotriacetonamine, velocity of catalysis of (McBAIN and BOLAM), 825.

9 IV

$C_9H_{10}O_4NBr$ Bromoaminoveratric acids (SIMONSEN and RAU), 789.

C_{10} Group.

$C_{10}H_{18}$ Hydrocarbon, from spinacene and sodium (CHAPMAN), 464.

10 II

$C_{11}H_8O_3$ 1-Hydrindone-2-carboxylic acid (ROBINSON and CRABTREE), 879.
$C_{11}H_8S_2$ β-Naphthylthiosulphoxylic acid, and its sodium salt (WHITE), 608.
$C_{10}H_9N$ β-Naphthylamine, sulphonation of (GREEN and VAKIL), 35.
$C_{10}H_{10}O_4$ Phenylsuccinic acids, optical activity of, and their esters (WREN), 210.
$C_{10}H_{11}Cl$ α-Chloro-*ar*-tetrahydronaphthalene (GREEN and ROWE), 971.
$C_{10}H_{12}O_4$ 5:6-Dimethoxy-*o*-toluic acid (PERKIN), 762.
$C_{10}H_{13}N$ *ar*-Tetrahydro-α-naphthylamine, preparation and reactions of (GREEN and ROWE), 955.
$C_{10}H_{15}N$ Aminobutylbenzene (REILLY and HICKINBOTTOM), 983.

10 III

$C_{10}H_4O_6N_4$ Dinitronaphth*iso*oxadiazole (GREEN and ROWE), 72.
$C_{10}H_5O_5N_3$ Nitronaphth*iso*oxadiazole (GREEN and ROWE), 72.
$C_{10}H_7ONa$ Sodium α-naphthoxide, relative activity of alkyl iodides with, in methyl alcohol (COX), 821.
Sodium α- and β-naphthoxides, relative activities of alkyl iodides with (COX), 666.
$C_{10}H_{10}O_5N_2$ 2:4-Dinitro-*ar*-tetrahydro-α-naphthol (GREEN and ROWE), 969.
$C_{10}H_{11}O_3N$ 2- and 4-Nitro-*ar*-tetrahydro-α-naphthol (GREEN and ROWE), 968.
$C_{10}H_{11}O_4N_3$ 2:4-Dinitro-*ar*-tetrahydro-α-naphthylamine, and its potassium salt (GREEN and ROWE), 962.
$C_{10}H_{12}O_3N_2$ 4-Nitro-*ar*-tetrahydro-α-naphthylamine, and its salts (GREEN and ROWE), 960.
$C_{10}H_{13}O_4N_4$ Semicarbazones of 5-nitro-2- and -6-methoxy-*m*-tolualdehydes (SIMONSEN), 778.
$C_{10}H_{11}O_4S$ *ar*-Tetrahydro-α-naphthol-4-sulphonic acid, and its sodium salt (GREEN and ROWE), 967.
$C_{10}H_{12}O_5N_2$ 5-Nitro-4-acetylaminoveratrole (SIMONSEN and RAU), 27.
$C_{10}H_{12}O_3N_2$ 2-Methoxy-*m*-tolualdehyde semicarbazone (SIMONSEN), 777.
$C_{10}H_{17}ON$ Camphoroxime, association of, in benzene (INNES), 432.

10 IV

$C_{10}H_3O_4N_2Cl$ 1-Chloro-2:4-dinitro-*ar*-tetrahydronaphthalene (GREEN and ROWE), 972.

$C_{10}H_{11}O_6NS$ 2-Nitro-*ar*-tetrahydro-α-naphthol-4-sulphonic acid (Gard and Rowe), 967.

$C_{10}H_{12}O_2NCl$ *p*-Chloromandeliminohydrin (Rule), 15.

$C_{10}H_{11}O_2NBr$ Bromoacetylaminoveratroles (Simonsen and Rau), 785, 790.

C_{11} Group.

$C_{11}H_{17}N$ *n*-Butyl-*o*- and -*p*-toluidines, and their salts (Reilly and Hickinbottom), 978.

$C_{11}H_{22}O_6$ Tetramethyl-γ-methylgalactoside (Cunningham), 599.

Tetramethyl-β-methylglucoside, preparation of (Haworth and Leitch), 194.

11 III

$C_{11}H_{12}O_7N_2$ 6-Nitro-5-acetylamino-3:4-dimethoxybenzoic acid (Simonsen and Rau), 25.

$C_{11}H_{13}O_4Br$ 6-Bromo-3:4-dimethoxy-β-phenylpropionic acid (Crabtree and Robinson), 871.

$C_{11}H_{13}O_5N$ Acetylamino-3:4-dimethoxybenzoic acids (Simonsen and Rau), 25.

$C_{11}H_{13}O_7N_5$ 2:3:5-Trinitro *p*-tolyl-*n*-butylnitrosoamine (Reilly and Hickinbottom), 993.

$C_{11}H_{13}O_6N_5$ 2:3:5-Trinitro-*p*-tolyl-*n*-butylnitroamine (Reilly and Hickinbottom), 993.

$C_{11}H_{14}O_5N_4$ 3:5-Dinitro-*p*-tolyl-*n*-butylnitrosoamine (Reilly and Hickinbottom), 991.

$C_{11}H_{14}O_6N_4$ 3:5-Dinitro-*p*-tolyl-*n*-butylnitroamine (Reilly and Hickinbottom), 992.

$C_{11}H_{15}O_3N_3$ 2-Nitro-*p*-tolyl-*n*-butylnitrosoamine (Reilly and Hickinbottom), 989.

$C_{11}H_{15}O_4N_3$ 3:5-Dinitro-*n*-butyl-*p*-toluidine (Reilly and Hickinbottom), 990.

$C_{11}H_{16}ON_2$ 5-Nitroso-*n*-butyl-*o*-toluidine, and its salts (Reilly and Hickinbottom), 982.

o- and *p*-Tolyl-*n*-butylnitrosoamine (Reilly and Hickinbottom), 979.

$C_{11}H_{16}O_2N_2$ Nitro-*n*-butyl-*p*-toluidines, and their salts (Reilly and Hickinbottom), 988.

11 IV

$C_{11}H_{12}O_5NBr$ Bromoacetylaminoveratric acids (Simonsen and Rau), 788.

C_{12} Group.

$C_{12}H_{10}N_2$ Azobenzene, association of, in benzene and alcohol (Innes), 430.

$C_{12}H_{22}O_{11}$ Lactose and Melibiose, constitution of (Haworth and Leitch), 188.

12 III

$C_{12}H_4O_6N_4$ Dinitronaphth*iso*xadiazole oxide (Green and Rowe), 71.

$C_{12}H_{13}O_8N_3$ 2:4-Dinitro-*ar*-tetrahydroaceto-α-naphthalide (Green and Rowe), 961.

$C_{12}H_{14}O_3N_2$ 4-Nitro-*ar*-tetrahydroaceto-α-naphthalide (Green and Rowe), 959.

$C_{12}H_{15}O_4N$ Cotarnine, sodium salt (Rakshit), 469.

$C_{12}H_{12}ON_2$ 1-Acetyl-*ar*-tetrahydro-1:4-naphthylenediamine (Green and Rowe), 959.

p-Hydroxybenzaldehydepiperylhydrazone (Weinhagen), 586.

Salicylaldehydepiperylhydrazone (Weinhagen), 585.

C_{13} Group.

$C_{13}H_{10}O$ Benzophenone, association of, in benzene (INNES), 430.

$C_{13}H_{18}N_2$ *p*-Aminoacetophenonepiperylhydrazone(+H_2O)(WEINHAGEN), 587.

$C_{13}H_{24}O_{11}$ Methyldigalactoside (CUNNINGHAM), 602.

13 III

$C_{13}H_{11}ON$ Formyldiphenylamine, association of, in benzene (INNES), 431.

$C_{13}H_{11}O_3N$ 2:4-Dihydroxybenzylideneaniline (SENIER and GALLAGHER), 33.

$C_{13}H_{15}O_2N_3$ Piperonalpiperylhydrazone (WEINHAGEN), 585.

$C_{13}H_{17}O_4N_3$ 3:5-Dinitroaceto-*n*-butyl-*p*-toluidide (REILLY and HICKINBOTTOM), 990.

$C_{13}H_{15}ON_3$ Anisaldehydepiperylhydrazone (WEINHAGEN), 586.

$C_{13}H_{19}O_3N_3$ Nitroaceto-*n*-butyl-*p*-toluidides (REILLY and HICKINBOTTOM), 989.

$C_{13}H_{19}ON$ Aceto-*n*-butyl-*o*- and -*p*-toluidides (REILLY and HICKINBOTTOM), 979.

13 IV

$C_{13}H_{10}O_2NBr$ 2:4-Dihydroxybenzylidenebromoanilines (SENIER and GALLAGHER), 34.

C_{14} Group.

$C_{14}H_{10}$ Phenanthrene, association of, in benzene and alcohol (INNES), 431.

14 II

$C_{14}H_8O_2$ 3:4-Phenanthraquinone (BARGER), 220.

$C_{14}H_{10}O_2$ Benzil, association of, in benzene and alcohol (INNES), 430.

3:4-Dihydroxyphenanthrene (*morphol*), synthesis of (BARGER), 218.

$C_{14}H_{23}N$ Di-*n*-butylaniline, and its salts (REILLY and HICKINBOTTOM), 99.

$C_{14}H_{16}N_2$ *p*-Phenylenedi-*n*-butyldiamine, and its hydrochloride (REILLY and HICKINBOTTOM), 107.

14 III

$C_{14}H_{13}O_2N$ γ-Benzilmonoxime, association of, in benzene (INNES), 432.

$C_{14}H_{13}O_3N$ 2:4-Dihydroxybenzylidene-*m*-toluidine (SENIER and GALLAGHER), 34.

$C_{14}H_{22}ON_2$ *p*-Nitrosodi-*n*-butylaniline, and its salts (REILLY and HICKINBOTTOM), 103.

C_{15} Group.

$C_{15}H_{12}O_3$ Morindone, constitution and reactions of (SIMONSEN), 766.

$C_{15}H_{10}O_6$ Santol (O'NEILL and PERKIN), 137.

$C_{15}H_{23}N$ Di-*n*-butyl-*p*-toluidine, and its salts (REILLY and HICKINBOTTOM), 980.

15 III

$C_{15}H_{10}O_{13}N_4$ 3:5-Dinitro-2-methoxyphenyl carbonate (POLLECOFF and ROBINSON), 649.

$C_{15}H_{12}O_9N_2$ 5-Nitro-2-methoxyphenyl carbonate (POLLECOFF and ROBINSON), 648.

$C_{15}H_{12}NCl$ Cinnamylidenechloroanilines (SENIER and GALLAGHER), 30.

$C_{15}H_{12}NBr$ Cinnamylidenebromoanilines (SENIER and GALLAGHER), 30.

$C_{15}H_{24}O_2N_2$ 2-Nitrodi-*n*-butyl-*p*-toluidine, and its hydrochloride (REILLY and HICKINBOTTOM), 994.

15 IV

$C_{15}H_{14}O_2NBr$ Benzoyl derivatives of bromoaminoveratroles (SIMONSEN and RAU), 786.

C_{16} Group.

$C_{16}H_{10}O_3$ 7-Hydroxy-4:3-indeno-1:2-benzoquinone (ROBINSON and CRABTREE), 879.

$C_{16}H_{12}O_3$ *r-Diphenylsuccinic anhydride, action of alcohols and amines on* (WREN and WILLIAMS), 832.

7-Methoxyphenylbenzopyrones (ROBINSON and TURNER), 875.

$C_{16}H_{12}O_5$ *iso*Brazilein, synthesis of salts of (CRABTREE and ROBINSON), 859.

Morindone methyl ether (SIMONSEN), 773.

$C_{16}H_{12}O_6$ Santal (O'NEILL and PERKIN), 136.

$C_{16}H_{14}O_4$ Diphenylsuccinic acids, optical activity of, and their esters (WREN), 210.

$C_{16}H_{16}O_3$ 2-Hydroxy-4-methoxyphenyl phenylethyl ketone (CRABTREE and ROBINSON), 870.

$C_{16}H_{32}O_2$ Palmitic acid, cerous salt (MORRELL), 116.

$C_{16}H_{34}O$ Hexadecyl alcohol, association of, in benzene and alcohol (INNES), 431.

16 III

$C_{16}H_{13}O_3Cl$ 7-Methoxy-2-phenyl-1:4-benzopyranol anhydrohydrochloride (+ $3H_2O$), (ROBINSON and TURNER), 877.

$C_{16}H_{14}O_2N_2$ Cinnamylidenenitrotoluidines (SENIER and GALLAGHER),

$C_{16}H_{15}ON$ Cinnamylidene-*p*-anisidine (SENIER and GALLAGHER), 31.

$C_{16}H_{12}O_4N_2$ 2:4-Dinitro-*ac*- and *-ar*-tetrahydro-α-naphthylamines (GREEN and ROWE), 972.

$C_{16}H_{18}O_2N_2$ Phenylacetiminohydrin (RULE), 11.

$C_{16}H_{18}O_4N_2$ Mandeliminohydrin, preparation of (RULE), 12.

16 IV

$C_{16}H_{19}O_3N_2S$ 4-*p*-Sulphobenzeneazo-*n*-butylaniline, sodium salt (REILLY and HICKINBOTTOM), 111.

C_{17} Group.

$C_{17}H_{14}O_3$ 3-Acetyl-2-phenyl-1:4-benzopyranol or 3-Benzoyl-2-methyl-1:4-benzopyranol (+ H_2O), and its anhydrohydrochloride (CHATTERJI and GHOSH), 446.

7-Hydroxy-3-benzyl-2-methyl-1:4-benzopyrone (CRABTREE and ROBINSON), 867.

Salicylidenebenzoylacetone (+ $\frac{1}{2}H_2O$), and its anhydrohydrochloride (CHATTERJI and GHOSH), 447.

$C_{17}H_{16}O_3$ Substance, from acetylacetone and salicylaldehyde (CHATTERJI and GHOSH), 448.

$C_{17}H_{16}O_4$ 2:4-Dimethoxydibenzoylmethane (ROBINSON and TURNER), 876.

Methyl hydrogen *r*-diphenylsuccinate, and its sodium salt (WREN and WILLIAMS), 837.

$C_{17}H_{17}N$ Cinnamylidene-*p*-xylidine (SENIER and GALLAGHER), 32.

17 III

$C_{17}H_{13}O_2N$ 2:4-Dihydroxybenzylidene-β-naphthylamine (SENIER and GALLAGHER), 37.

$C_{17}H_{17}ON$ Cinnamylidene-*p*-phenetidine (SENIER and GALLAGHER), 32.

$C_{17}H_{19}O_3N$ Morphine, calcium salt (RAKSHIT), 470.

$C_{17}H_{20}O_3N_4$ Osazone of sugar from morindin (SIMONSEN), 774.

$C_{17}H_{23}O_3N$ 2:4-Dihydroxybenzylidenecamphylamine (SENIER and GALLAGHER), 35.

17 IV

$C_{17}H_{21}O_3N_2S$ 4-Methyl-*N*-*n*-butyldiazoaminobenzene-4'-sulphonic acid, and its salts (REILLY and HICKINBOTTOM), 984.

C_{18} Group.

$C_{16}H_{16}O_5$ 7-Hydroxy-3-benzyl-2-methyl-1:4-benzopyrone methyl ether (Crabtree and Robinson), 867.

$C_{18}H_{16}O_5$ Morindone trimethyl ether (Simonsen), 772.

$C_{18}H_{18}O_4$ Ethyl hydrogen *r*- and *meso*-diphenylsuccinates, and their metallic salts (Wren and Williams), 835.

$C_{18}H_{20}O_5$ 2-Hydroxy-4-methoxyphenyl 3:4-dimethoxyphenylethyl ketone (Crabtree and Robinson), 871.

$C_{18}H_{32}O_4$ Menthyl hydrogen terephthalate (Cohen and de Pennington), 64.

$C_{18}H_{30}O_2$ Linolenic acid, cerous salt (Morrell), 119.

$C_{18}H_{32}O_2$ *α*- and *β*-Elaeostearic acids, cerous and lead salts (Morrell), 117. Linoleic acid, cerous salt (Morrell), 117.

$C_{18}H_{34}O_2$ Elaidic acid, cerous salt (Morrell), 117. Oleic acid, and its potassium salt, molecular condition of, in alcoholic solution (Laing), 435; cerous salt (Morrell), 117.

$C_{18}H_{36}O_2$ Stearic acid, cerous salt (Morrell), 112.

18 III

$C_{15}H_{13}O_2N_3$ Formylmethylaminobenzene-4-azo-*β*-naphthol (Morgan and Grist), 692.

$C_{18}H_{18}O_3N_2$ Cinnamylidenenitro-ψ-cumidine (Senier and Gallagher), 32.

$C_{18}H_{19}O_4N$ α-Imino-*ββ*-diphenyldipropionic acid (Senter, Drew, and Martin), 161.

$C_{15}H_{21}ON$ Benzo-*n*-butyl-*p*-toluidide (Reilly and Hickinbottom), 979.

$C_{18}H_{21}O_3N$ Codeine, potassium and sodium salts of (Rakshit), 466.

$C_{18}H_{23}O_4N$ Menthyl hydrogen nitroterephthalate (Cohen and de Pennington), 64.

C_{19} Group.

$C_{19}H_{16}O_4$ 7-Acetoxy-3-benzyl-2-methyl-1:4-benzopyrone (Crabtree and Robinson), 867.

$C_{19}H_{28}O_4$ Methyl menthyl terephthalate (Cohen and de Pennington), 63.

$C_{19}H_{34}O_{11}$ Hexamethyl methyllactoside (Haworth and Leitch), 195.

19 III

$C_{19}H_{25}O_6N$ *o*-Nitromethyl menthyl terephthalate (Cohen and de Pennington), 66.

C_{20} Group.

$C_{20}H_{14}O_4$ 3′:4′-Methylenedioxy-2-phenyl-2:3-dihydro-1:4-*α*-naphthapyrone (Crabtree and Robinson), 865.

$C_{20}H_{14}O_7$ Diacetylmorindone methyl ether (Simonsen), 774.

$C_{20}H_{20}O_4$ Berberidene (Perkin), 761.

$C_{20}H_{20}O_5$ Ketodihydroberberidene (Perkin), 763.

$C_{20}H_{28}O_4$ Ethyl menthyl terephthalate (Cohen and de Pennington), 63.

$C_{20}H_{30}O_2$ Abietic acid, cerous salt (Morrell), 120.

$C_{20}H_{38}O_{11}$ Heptamethyl methyllactoside (Haworth and Leitch), 195. Octamethyl digalactose (Cunningham), 601.

20 III

$C_{20}H_{13}O_2N$ *N*-Phenoxyacetylcarbazole (Copisarow), 818.

$C_{20}H_{17}O_2N_3$ Acetylmethylaminobenzene-4-azo-*β*-naphthol (Morgan and Grist), 694.

$C_{20}H_{17}O_5N$ Oxyberberine, preparation of (Perkin), 737. *α*- and *iso*-Oxy*epi*berberines (Perkin), 518.

$C_{20}H_{19}O_4N$ Dihydroanhydro-*n*- and -*epi*-berberines, and their salts (PERKIN 506, 737.

$C_{20}H_{19}O_5N$ *epi*Berberine, and its salts (PERKIN), 492.

$C_{20}H_{21}O_4N$ Tetrahydroanhydro*epi*berberine, and its salts PERKIN), 510.

$C_{20}H_{27}O_4N$ Nitroethyl menthyl terephthalates (COHEN and DE P NINGTON), 66.

20 IV

$C_{20}H_{27}O_3N_2S$ 4-*p*-Sulphobenzeneazodi-*n*-butylaniline, sodium and po sium salts (REILLY and HICKINBOTTOM), 109.

C_{21} Group.

$C_{21}H_{16}O_8$ Acetylmorindone (SIMONSEN), 770.

21 III

$C_{21}H_{14}O_3N_2$ Hydroxymethoxyphenanthraphenazines (POLLECOFF ; ROBINSON), 651.

$C_{21}H_{20}O_7N_2$ 5:5'-Dinitro-2:2'- and -4:4'-dimethoxydi-3-*m*-methylst ketones (SIMONSEN), 778.

$C_{21}H_{21}O_4N$ Anhydromethylberberines, and their salts (PERKIN), 746, 75

$C_{21}H_{21}O_5N$ *O*-Methyl*epi*berberine (PERKIN), 520.

$C_{21}H_{21}O_7N$, Trioxyanhydromethylberberine (PERKIN), 747.

$C_{21}H_{23}O_5N$ Hydroxy*iso*anhydrodihydromethylberberines (PERKIN),

$C_{21}H_{23}O_5N_3$ Semicarbazone of ketodihydroberberidene (PERKIN), 76

$C_{21}H_{25}ON$ *N*-*n*-Nonoylcarbazole (COPISAROW), 818.

$C_{21}H_{25}O_4N$ Dihydromethyl*iso*tetrahydroanhydroberberine, and its s (PERKIN), 759.

21 IV

$C_{21}H_{22}O_4NCl$ ψ-Methylberberinium chloride ($+H_2O$) (PERKIN), 750.

$C_{21}H_{22}O_4NI$ ψ-Berberinium iodide (PERKIN), 751.

$C_{21}H_{24}O_4NI$ *N*-Methyl*iso*tetrahydroanhydroberberine hydriodide (PE KIN), 748.

C_{22} Group.

$C_{22}H_{18}O_3$ 7-Methoxy-2:4-diphenyl-1:4-benzopyranol, and its salts (ROBI SON and TURNER), 877.

22 III

$C_{22}H_{19}O_3N$ *r*-Diphenylsuccinanilic acid (WREN and WILLIAMS), 837.

$C_{22}H_{25}O_5N$ *O*-Ethyl*epi*berberine (PERKIN), 521.

$C_{22}H_{23}O_7N$ Narcotine, potassium and sodium salts (RAKSHIT), 467.

$C_{22}H_{27}O_4N$ Dihydrodimethyl*iso*tetrahydroanhydroberberine, and i salts (PERKIN), 760.

$C_{22}H_{31}O_6N$ Nitrobutyl menthyl terephthalates (COHEN and DE PE NINGTON), 66.

C_{23} Group.

$C_{23}H_{19}O_2N$ Diphenylsuccino-*p*-tolil (WREN and WILLIAMS), 839.

$C_{23}H_{21}O_3N$ *meso*- and *r*-Diphenylsuccino-*p*-toluidic acid, and the silver s of the latter (WREN and WILLIAMS), 838.

$C_{23}H_{21}O_2N_3$ Phenylhydrazone of substance $C_{17}H_{16}O_3$ (CHATTERJI an GHOSH), 449.

$C_{23}H_{23}O_4N$ Anhydro*epi*berberineacetone (PERKIN), 521.

$C_{23}H_{25}O_4N$ Acetoxy*iso*anhydrodihydromethylberberines (PERKIN), 75

C_{24} Group.

$C_{24}H_{20}N_2$ Dicinnamylidene-*p*-phenylenediamine (SENIER and GA LAGHER), 32.

$C_{21}H_{22}O_5$ Substance, from acetylacetone and salicylaldehyde (CHATTERJI and GHOSH), 449.
$C_{24}H_{24}O_7$ Deoxy-*n*- and -*iso*-santalin (O'NEILL and PERKIN), 130, 133.
$C_{24}H_{24}O_8$ *n*- and *iso*-Santalin (O'NEILL and PERKIN), 127, 131.
$C_{24}H_{34}O_4$ *cyclo*Hexyl menthyl terephthalate (COHEN and DE PENNINGTON), 63.

24 III

$C_{31}H_{25}O_3N$ Methyl *r*-diphenylsuccino-*p*-toluidate (WREN and WILLIAMS), 839.
$C_{24}H_{39}ON_3$ 4-Di-*n*-butylaminobenzeneazo-β-naphthol (REILLY and HICKINBOTTOM), 108.

C_{25} Group.

$C_{25}H_{22}O_7$ Santalone (O'NEILL and PERKIN), 133.
$C_{25}H_{44}O_{21}$ Methyltetragalactoside and Methyltetraglucoside (CUNNINGHAM), 606.

25 III

$C_{25}H_{16}ON_2$ *N*-Carbonylcarbazole (COPISAROW), 819.

C_{26} Group.

$C_{26}H_{22}N_2$ Dicinnamylidene-*o*-naphthylenediamine (SENIER and GALLAGHER), 33.
$C_{26}H_{40}O_4$ Octyl menthyl terephthalate (COHEN and DE PENNINGTON), 63.

26 III

$C_{28}H_{16}O_2N_2$ *N*-Oxalylcarbazole (COPISAROW), 819.

C_{27} Group.

$C_{27}H_{20}O_2$ Substance ($+\frac{1}{2}H_2O$), from benzoylacetone and salicylaldehyde (CHATTERJI and GHOSH), 446.

C_{28} Group.

$C_{28}H_{18}O_4$ α-Naphtholphthalein, preparation of (WERNER), 20.
$C_{28}H_{26}O_6$ Piperonylidene derivative of 3′:4′-methylenedioxy-2-phenyl-2:3-dihydro-1:4-α-naphthapyrone (CRABTREE and ROBINSON), 865.
$C_{28}H_{21}N$ Tetraphenylpyrrole, synthesis of (G. M. and R. ROBINSON), 639.
$C_{28}H_{42}O_4$ Menthyl terephthalate (COHEN and DE PENNINGTON), 64.

28 III

$C_{28}H_{39}ON$ *N*-Palmitylcarbazole (COPISAROW), 819.
$C_{28}H_{41}O_6N$ Menthyl nitroterephthalate (COHEN and DE PENNINGTON), 64.

C_{29} Group.

$C_{29}H_{48}$ Spinacene, analysis and reactions of, and its hexahydro-bromide and -chloride (CHAPMAN), 458.

29 III

$C_{29}H_{40}O_4N_2$ *iso*Emetine ($+H_2O$), and its salts (PYMAN), 226.

C_{30} Group.

$C_{30}H_{24}N_2$ Dicinnamylidenebenzidine (SENIER and GALLAGHER), 33.

30 III

$C_{30}H_{42}O_4N_2$ *N*-Methyl*iso*emetine (PYMAN), 228.

C_{32} Group.

$C_{32}H_{30}O_{12}$ Acetyl-*n*- and -*iso*-santalin (O'NEILL and PERKIN), 128.
$C_{32}H_{32}O_{11}$ Acetyldeoxy-*n*- and -*iso*-santalin (O'NEILL and PERKIN), 130, 134.

32 III

$C_{32}H_{48}O_4N_2$ *N*-Methyl*iso*emetinemethine, and its salts (PYMAN), 229.

32 IV

$C_{32}H_{46}O_4N_2Cl_2$ *N*-Methylemetine methochlorides (PYMAN), 233.
$C_{32}H_{46}O_4N_2I_2$ *N*-Methyl-*n*- and -*iso*-emetine methiodides (PYMAN), 233.

C_{35} Group.

$C_{35}H_{39}O_5N_5$ Ergotinine, supposed formation of ergotoxine ethyl ester fro (BARGER and EWINS), 235.
$C_{35}H_{41}O_6N_5$ Ergotoxine, formation of the phosphate of, from ergotinin (BARGER and EWINS), 235.

C_{36} Group.

$C_{36}H_{32}O_8$ Benzoylmorindone (SIMONSEN), 771.

C_{37} Group.

$C_{37}H_{64}O_{31}$ Methyltrimaltoside (+ H_2O) (CUNNINGHAM), 607

ERRATA.

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Page	Line	
243	15*	*for* "water" *read* "oxygen."
774	15*	,, "sodium hypochlorite" *read* "potassium permanganate."

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31	19	,, "$C_{16}H_{15}O_3N_3$" *read* "$C_{16}H_{14}O_3N_2$."
100	3*	,, "4-*di*-n-*butylaminoazobenzene*-4′-*sulphonic acid*" read "4-*sulphobenzeneazodi*-n-*butylaniline*."
100	1	,, "4-n-*butylaminoazobenzene*-4′-*sulphonic acid*" read "4-p-*sulph benzeneazo*-n-*butylaniline*."
122	16*	,, "$Ce_3(C_{18}H_{31}O_3.O)_3$" *read* "$Ce(C_{18}H_{31}O_3.O_2)_3$."
123	4	,, "(Average temperature)" *read* "(Room temperature)."
540	4	,, "+ 50·6°" *read* "− 50·6°."
774	6	,, "$C_{20}H_{16}O_7$" *read* "$C_{20}H_{14}O_7$."

* From bottom.